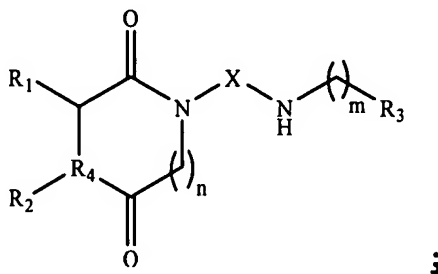


Listing of Claims

1. (currently amended) A compound ~~of formula I, a stereochemical isomer of the compound, or a hydrate, solvate, or pharmaceutically acceptable salt of the compound or isomer, wherein:~~

the compound corresponds in structure to formula I:



~~one of their stereochemically isomer forms or a pharmaceutically acceptable salt thereof, wherein:~~

R₄ is selected from the group consisting of N and S;

if R₄ is S, then R₁ is H, and R₂ is absent;

if R₄ is N, then R₁ and R₂ are H or are methylene groups bound together forming with the heterocyclic ring a 5- or 6- membered ring; if R₄=S then R₁ is H and R₂ is absent; R₄ is selected from the group consisting of N and S;

n being an integer from 0 to is zero or 1;

X is selected from the group consisting of C₂-C₁₀-alkyl~~ene~~, C₂-C₁₀-alkenyl~~ene~~, and -CH₂-Y-CH₂-; ~~wherein~~

Y is phenyl;

~~m being an integer from~~ m is 1 [[to]] or 2;

R₃ is selected from the group consisting of chroman-2-yl, 2-quinolyl, and phenoxy -O-phenyl, wherein:

the quinolyl, the aromatic ring of the chromanyl ~~moiety, the quinolyl, and [[or]]~~ the phenyl ring of the phenoxy are residue is optionally substituted [[by]] with one or more groups chosen substituents independently selected from the group consisting of C₁-C₆-alkoxy, C₁-C₆-alkyl, halogen, C₂-C₆-alkenyl, halo-(C₁-C₆)-alkyl, halo-(C₁-C₆)-

alkoxy, phenyl, phenyl(C₁-C₆)-alkyl, phenoxy, C₁-C₆-alkylcarbonyl, phenylcarbonyl, phenyl(C₁-C₆)alkylcarbonyl, C₁-C₆-alkoxycarbonyl, phenyl(C₁-C₆)alkoxycarbonyl, C₁-C₆-alkyl-carbonylamino, hydroxy, cyano, nitro, amino, N-(C₁-C₆)-alkylamino, N,N-(C₁-C₆)-dialkylamino, carboxy, sulfo, sulfamoyl, sulfonylamino, (C₁-C₆)alkylaminosulfonyl, and [[or]] (C₁-C₆)alkylsulfonylamino, wherein:

the C₁-C₆-alkyl portion of any of the alkyl-comprising substituents is optionally substituted with a substituent independently selected from the group consisting of hydroxy and amino; or wherein

the phenyl ring of the phenoxy is substituted by two neighbouring residues, which together with the phenyl ring to which they are attached form tetrahydronaphthyl; ~~wherein each alkyl is optionally substituted with hydroxy or amino;~~ provided that the compound is not

2-[4-[(chroman-2-yl)methylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-c]imidazole,
3-[4-[(chroman-2-yl)methylamino]butyl]-2,4-dioxothiazolidine,
3-[5-[(chroman-2-yl)methylamino]pentyl]-2,4-dioxothiazolidine,
3-[6-[(chroman-2-yl)methylamino]hexyl]-2,4-dioxothiazolidine,
2-[4-[2-(phenoxy)ethylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-c]imidazole, [[or]]
3-[4-[2-(phenoxy)ethylamino]butyl]-2,4-dioxothiazolidine, or and is not
3-[3-[(chroman-2-yl)methylamino]propyl]-2,4-dioxoimidazolidine dioxoimidazoidine.

2. (currently amended) ~~Compound~~ The compound, isomer, hydrate, solvate, or salt according to claim 1, wherein R₃ is selected from the group consisting of chroman-2-yl, 2-quinolyl, and ~~phenoxy -O-phenyl~~, wherein:

the phenyl ring of the phenoxy residue is optionally substituted ~~by a group chosen with a substituent selected~~ from the group consisting of C₁-C₆-alkoxy, C₁-C₆-alkyl, [[or]] and halogen; [[;]]

3. (currently amended) ~~Compound~~ The compound, isomer, hydrate, solvate, or salt according to claim 1 [[or 2]], wherein:
m is 1; and

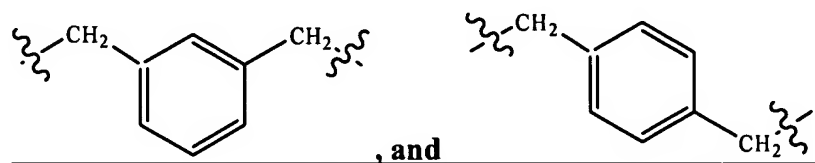
R₃ is chroman-2-yl.

4. (currently amended) ~~Compound~~ The compound, isomer, hydrate, solvate, or salt according to claim 3, wherein:

R₁ and R₂ are methylene groups bound together forming with the heterocyclic ring a 5- or 6- membered ring; and

R₄ is N.

5. (currently amended) ~~Compound~~ The compound, isomer, hydrate, solvate, or salt according to ~~any of claims~~ claim 3 [[to 4]], wherein X is selected from the group consisting of C₂-C₁₀-alkylene, (E)-2-butenylene, ~~3-methylbenzyl or 4-methylbenzyl~~.



6. (currently amended) ~~Compound~~ The compound, isomer, hydrate, solvate, or salt according to claim 3, wherein:

R₁ is H₂ [[,]]

R₂ is absent; and

R₄ is S.

7. (currently amended) ~~Compound~~ The compound, isomer, hydrate, solvate, or salt according to claim 6, wherein:

n is zero; [[0]] and

X is C₂-C₁₀-alkylene.

8. (currently amended) ~~Compound~~ The compound, isomer, hydrate, solvate, or salt according to claim 1 [[or 2]], wherein:

m is [[=]] 2; and

R₃ is phenoxy -~~O-phenyl~~, wherein the phenyl ring of the phenoxy: residue

is optionally substituted [[by]] with one or more ~~groups-chosen~~ substituents

independently selected from the group consisting of C₁-C₆-alkoxy, C₁-C₆-alkyl, halogen, C₂-C₆-alkenyl, halo-(C₁-C₆)-alkyl, halo-(C₁-C₆)-alkoxy, phenyl, phenyl(C₁-C₆)-alkyl, phenoxy, C₁-C₆-alkylcarbonyl, phenylcarbonyl, phenyl(C₁-C₆)alkylcarbonyl, C₁-C₆-alkoxycarbonyl, phenyl(C₁-C₆)alkoxycarbonyl, C₁-C₆-alkylcarbonylamino, hydroxy, cyano, nitro, amino, N-(C₁-C₆)-alkylamino, N,N-(C₁-C₆)-dialkylamino, carboxy, sulfo, sulfamoyl, sulfonylamino, (C₁-C₆)alkylaminosulfonyl, and [[or]] (C₁-C₆)alkylsulfonylamino; or ~~wherein the phenyl ring~~

is substituted by two neighbouring residues, which together with the phenyl ring to which they are attached form tetrahydronaphthyl.

9. (currently amended) ~~Compound~~ The compound, isomer, hydrate, solvate, or salt according to claim 8, wherein:

R₃ is phenoxy, wherein the phenyl [[group]] ring of the phenoxy:

is optionally substituted [[by]] with one or more ~~groups-chosen substituents~~ independently selected from the group consisting of phenyl, C₁-C₆-alkoxycarbonyl, C₁-C₆-alkylcarbonylamino, C₁-C₆-alkoxy, C₁-C₆-alkyl, halo-(C₁-C₆)-alkyl, and [[or]] halogen, or ~~wherein the phenyl group~~

is substituted by two neighbouring neighbouring residues, which together with the phenyl group to which they are attached form tetrahydronaphthyl.

10. (currently amended) ~~Compound~~ The compound, isomer, hydrate, solvate, or salt according to claim 9, wherein the phenyl ~~residue~~ ring of the phenoxy is optionally substituted [[by]] with one or more ~~groups-chosen substituents independently selected~~ from the group consisting of methoxy, ethoxy, propoxy, isopropoxy, ethyl, propyl, isopropyl, bromide, trifluoromethyl, ~~methylamide~~ methylamido, and [[or]] ethoxycarbonyl.

11. (currently amended) ~~Compound~~ The compound, isomer, hydrate, solvate, or salt according to ~~any of claims~~ claim 8 [[to 10]], wherein R₃ is phenoxy, wherein:
the phenyl ~~group~~ ring of the phenoxy is substituted in ortho- and/or meta- position.

12. (currently amended) ~~Compound~~ The compound, isomer, hydrate, solvate, or salt according to ~~any of claims~~ claim 8 [[to 11]], wherein:

R₁ and R₂ are methylene groups bound together forming with the heterocyclic ring a 5- or 6- membered ring; and

R₄ is N.

13. (currently amended) ~~Compound~~ The compound, isomer, hydrate, solvate, or salt according to ~~any of claims~~ claim 8 [[to 12]], wherein:

n is [[0]] zero; and

X is C₂-C₁₀-alkylene.

14. (currently amended) ~~Compound~~ The compound, isomer, hydrate, solvate, or salt according to ~~any of claims~~ claim 8 [[to 11]], wherein:

R₁ is H; [[and]]

R₂ is absent; and

R₄ is S.

15. (currently amended) ~~Compound~~ The compound, isomer, hydrate, solvate, or salt according to claim 14, wherein:

n is [[0]] zero; and

X is C₂-C₁₀-alkylene.

16. (currently amended) ~~Compound~~ The compound, isomer, hydrate, solvate, or salt according to ~~claims~~ claim 1 [[or 2]], wherein:

m is 1; and

R₃ is 2-quinolyl.

17. (currently amended) ~~Compound~~ The compound, isomer, hydrate, solvate, or salt according to claim 16, wherein:

R₁ and R₂ are methylene groups bound together forming with the heterocyclic ring a 5- or 6- membered ring; and

R₄ is N.

18. (currently amended) ~~Compound~~ The compound, isomer, hydrate, solvate, or salt according to ~~any of claims~~ claim 16 ~~[[to 17]]~~, wherein:

n is zero ~~[[0]]~~; and

X is C₂-C₁₀-alkylene.

19. (currently amended) ~~Compound~~ The compound, isomer, hydrate, solvate, or salt according to claim 1, wherein the compound is selected from the group consisting of:

~~(a) 2-[4-[(Chroman-2(R)-yl)methylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-c]imidazole;~~
[[b)]]

2-[4-[(Chroman-2-yl)methylamino]butyl]-1,3-dioxoperhydroimidazo[1,5-a]pyridine;

[[c)]]

2-[4-[(Chroman-2-yl)methylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-a]pyrazine;

[[d)]]

2-[5-[(Chroman-2-yl)methylamino]pentyl]-1,3-dioxoperhydropyrrolo[1,2-c]imidazole;

[[e)]]

2-[6-[(Chroman-2-yl)methylamino]hexyl]-1,3-dioxoperhydropyrrolo[1,2-c]imidazole;

[[f)]]

2-[3-[(Chroman-2-yl)methylamino]propyl]-1,3-dioxoperhydropyrrolo[1,2-c]imidazole;

[[g)]]

3-[8-[(Chroman-2-yl)methylamino]octyl]-2,4-dioxothiazolidine;

~~(h) 2-[4-[(Chroman-2(S)-yl)methylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-c]imidazole;~~
2-[4-[(Chroman-2-yl)methylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-c]

imidazole;

[[i)]]

2-[8-[(Chroman-2-yl)methylamino]octyl]-1,3-dioxoperhydropyrrolo[1,2-c]imidazole;

[[j)]]

2-[3-[[[(Chroman-2-yl)methylamino]methyl]benzyl]-1,3-dioxoperhydropyrrolo[1,2-c]

imidazole;

[[k)]]

2-[4-[[~~(Chroman-2-yl)methylamino~~]methyl]benzyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]
imidazole;

~~(1) (E) 2-[4-[(Chroman-2-yl)methylamino]but-2-enyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]
imidazole;~~

2-[4-[(Chroman-2-yl)methylamino]but-2-enyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]
imidazole;

[[**(m)**]]

2-[4-[2-(*o*-Methoxyphenoxy)ethylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]
imidazole;

[[**(n)**]]

2-[4-[2-(*m*-Methoxyphenoxy)ethylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]
imidazole;

[[**(o)**]]

2-[4-[2-(*o*-Bromophenoxy)ethylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]imidazole;

[[**(p)**]]

2-[4-[2-(*m*-Bromophenoxy)ethylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]imidazole;

[[**(q)**]]

2-[4-[2-(*o*-Ethylphenoxy)ethylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]imidazole;

[[**(r)**]]

2-[4-[2-(*m*-Ethylphenoxy)ethylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]imidazole;

[[**(s)**]]

2-[4-[2-(*o*-Isopropylphenoxy)ethylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]
imidazole;

[[**(t)**]]

2-[4-[(2-quinolyl)methylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]imidazole;

[[**(u)**]]

2-[4-[2-(*o*-Ethoxyphenoxy)ethylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]imidazole;

[[**(v)**]]

2-[4-[2-(*o*-Isopropoxyphenoxy)ethylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]

imidazole;

[[w]]

2-[4-[2-[m-(Trifluoromethyl)phenoxy]ethylamino]butyl]-1,3-dioxoperhydropyrrolo

[1,2-c] imidazole;

[[x]]

2-[4-[2-(1,1'-Biphenyl-2-yloxy)ethylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-c]

imidazole;

[[y]]

2-[4-[2-[o-(Acetylamino)phenoxy]ethylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-c]

imidazole;

[[z]]

2-[4-[2-[m-(Acetylamino)phenoxy]ethylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-c]

imidazole;

[[aa]]

2-[4-[2-[o-(Ethoxycarbonyl)phenoxy]ethylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-c]

imidazole;

[[bb]]

2-[4-[2-(5,6,7,8-Tetrahydronaphth-1-yloxy)ethylamino]butyl]-1,3-dioxoperhydropyrrolo

[1,2-c]imidazole;

[[cc]]

2-[4-[2-(2,3-Dimethylphenoxy)ethylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-c]

imidazole;

[[dd]]

2-[4-[(Chroman-2-yl)methylamino]butyl]-1,4-dioxoperhydropyrido[1,2-a]pyrazine;

~~(ee) (Z)-2-[4-[(Chroman-2-yl)methylamino]but-2-enyl]-1,4-dioxoperhydropyrrolo[1,2-c]~~

~~imidazole;~~

2-[4-[(Chroman-2-yl)methylamino]but-2-enyl]-1,4-dioxoperhydropyrrolo[1,2-c]

imidazole;

[[ff]]

3-[4-[2-(*o*-Ethoxyphenoxy)ethylamino]butyl]-2,4-dioxothiazolidine;

[[**(gg)**]]

3-[6-[2-(*o*-Ethoxyphenoxy)ethylamino]hexyl]-2,4-dioxothiazolidine;

[[**(hh)**]]

3-[8-[2-(*o*-Ethoxyphenoxy)ethylamino]octyl]-2,4-dioxothiazolidine;

[[**(ii)**]]

2-[4-[2-(*o*-Ethoxyphenoxy)ethylamino]butyl]-1,3-dioxoperhydroimidazo[1,5-*a*]pyridine;

[[**(jj)**]]

2-[6-[2-(*o*-Ethoxyphenoxy)ethylamino]hexyl]-1,3-dioxoperhydroimidazo[1,5-*a*]pyridine;

[[**(kk)**]]

2-[4-[(2-Quinolyl)methylamino]butyl]-1,3-dioxoperhydroimidazo[1,5-*a*]pyridine; and

[[**(ll)**]]

2-[6-[(2-Quinolyl)methylamino]hexyl]-1,3-dioxoperhydroimidazo[1,5-*a*]pyridine;

~~a pharmaceutically acceptable salt or one of their stereochemically isomer forms.~~

20. (currently amended) ~~Pharmaceutical~~ A pharmaceutical composition which comprises;

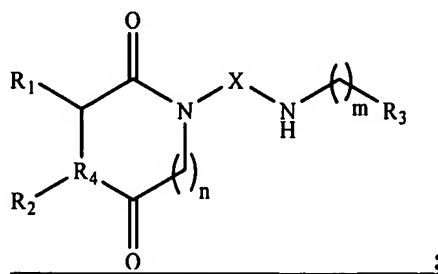
a therapeutically effective amount of a compound, a stereochemical isomer of the compound, or a hydrate, solvate, or pharmaceutically acceptable salt of the compound or isomer, wherein the compound is selected from the group of compounds recited in claim 1; as claimed in any of claims 1 to 19 and [[,]]

one or more pharmaceutically acceptable carriers.

21. (currently amended) [[Use]] A use of a compound, a stereochemical isomer of the compound, or a hydrate, solvate, or pharmaceutically acceptable salt of the compound or isomer of formula I according to any of claims 1 to 19, wherein the disclaimer to 3-[3-{(chroman-2-yl)methylamino}propyl]-2,4-dioxoimidazolidine does not apply, for the preparation of a medicament for the treatment and/or prophylaxis of a condition selected from the group consisting of Parkinson Disease, cerebral damage by thromboembolic ictus, craneoencephalic traumatism, depression, migraine, pain, psychosis, anxiety disorders,

aggressive disorders, and [[or]] urinary tract disorders, wherein:

the compound corresponds in structure to formula I:



R₄ is selected from the group consisting of N and S;

if R₄ is S, then R₁ is H, and R₂ is absent;

if R₄ is N, then R₁ and R₂ are H or are methylene groups bound together forming with the heterocyclic ring a 5- or 6- membered ring;

n is zero or 1;

X is selected from the group consisting of C₂-C₁₀-alkylene, C₂-C₁₀-alkenylyne, and -CH₂-Y-CH₂-;

Y is phenyl;

m is 1 or 2;

R₃ is selected from the group consisting of chroman-2-yl, 2-quinolyl, and phenoxy, wherein:

the quinolyl, the aromatic ring of the chromanyl, and the phenyl ring of the phenoxy are optionally substituted with one or more substituents independently selected from the group consisting of C₁-C₆-alkoxy, C₁-C₆-alkyl, halogen, C₂-C₆-alkenyl, halo-(C₁-C₆)-alkyl, halo-(C₁-C₆)-alkoxy, phenyl, phenyl(C₁-C₆)-alkyl, phenoxy, C₁-C₆-alkylcarbonyl, phenylcarbonyl, phenyl(C₁-C₆)alkylcarbonyl, C₁-C₆-alkoxycarbonyl, phenyl(C₁-C₆)alkoxycarbonyl, C₁-C₆-alkyl-carbonylamino, hydroxy, cyano, nitro, amino, N-(C₁-C₆)-alkylamino, N,N-(C₁-C₆)-dialkylamino, carboxy, sulfo, sulfamoyl, sulfonylamino, (C₁-C₆)alkylaminosulfonyl, and (C₁-C₆)alkylsulfonylamino, wherein:

the C₁-C₆-alkyl portion of any of the alkyl-comprising

substituents is optionally substituted with a substituent independently
selected from the group consisting of hydroxy and amino; or
the phenyl ring of the phenoxy is substituted by two neighbouring residues,
which together with the phenyl ring to which they are attached form
tetrahydronaphthyl;
provided that the compound is not
2-[4-[(chroman-2-yl)methylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-c]imidazole,
3-[4-[(chroman-2-yl)methylamino]butyl]-2,4-dioxothiazolidine,
3-[5-[(chroman-2-yl)methylamino]pentyl]-2,4-dioxothiazolidine,
3-[6-[(chroman-2-yl)methylamino]hexyl]-2,4-dioxothiazolidine,
2-[4-[2-(phenoxy)ethylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-c]imidazole, or
3-[4-[2-(phenoxy)ethylamino]butyl]-2,4-dioxothiazolidine.

22. (new) The use according to claim 21, wherein the compound is selected from the group consisting of:

2-[4-[(Chroman-2-yl)methylamino]butyl]-1,3-dioxoperhydroimidazo[1,5-a]pyridine;
2-[4-[(Chroman-2-yl)methylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-a]pyrazine;
2-[5-[(Chroman-2-yl)methylamino]pentyl]-1,3-dioxoperhydropyrrolo[1,2-c]imidazole;
2-[6-[(Chroman-2-yl)methylamino]hexyl]-1,3-dioxoperhydropyrrolo[1,2-c]imidazole;
2-[3-[(Chroman-2-yl)methylamino]propyl]-1,3-dioxoperhydropyrrolo[1,2-c]imidazole;
3-[8-[(Chroman-2-yl)methylamino]octyl]-2,4-dioxothiazolidine;
2-[4-[(Chroman-2-yl)methylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-c]imidazole;
2-[8-[(Chroman-2-yl)methylamino]octyl]-1,3-dioxoperhydropyrrolo[1,2-c]imidazole;
2-[3-[[[(Chroman-2-yl)methylamino]methyl]benzyl]-1,3-dioxoperhydropyrrolo[1,2-c]
imidazole;
2-[4-[[[(Chroman-2-yl)methylamino]methyl]benzyl]-1,3-dioxoperhydropyrrolo[1,2-c]
imidazole;
2-[4-[(Chroman-2-yl)methylamino]but-2-enyl]-1,3-dioxoperhydropyrrolo[1,2-c]
imidazole;

2-[4-[2-(*o*-Methoxyphenoxy)ethylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]
imidazole;

2-[4-[2-(*m*-Methoxyphenoxy)ethylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]
imidazole;

2-[4-[2-(*o*-Bromophenoxy)ethylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]imidazole;

2-[4-[2-(*m*-Bromophenoxy)ethylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]imidazole;

2-[4-[2-(*o*-Ethylphenoxy)ethylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]imidazole;

2-[4-[2-(*m*-Ethylphenoxy)ethylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]imidazole;

2-[4-[2-(*o*-Isopropylphenoxy)ethylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]
imidazole;

2-[4-[(2-quinolyl)methylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]imidazole;

2-[4-[2-(*o*-Ethoxyphenoxy)ethylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]imidazole;

2-[4-[2-(*o*-Isopropoxyphenoxy)ethylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]
imidazole;

2-[4-[2-[*m*-(Trifluoromethyl)phenoxy]ethylamino]butyl]-1,3-dioxoperhydropyrrolo
[1,2-*c*]imidazole;

2-[4-[2-(1,1'-Biphenyl-2-yloxy)ethylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]
imidazole;

2-[4-[2-[*o*-(Acetylamino)phenoxy]ethylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]
imidazole;

2-[4-[2-[*m*-(Acetylamino)phenoxy]ethylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]
imidazole;

2-[4-[2-[*o*-(Ethoxycarbonyl)phenoxy]ethylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]
imidazole;

2-[4-[2-(5,6,7,8-Tetrahydronaphth-1-yloxy)ethylamino]butyl]-1,3-dioxoperhydropyrrolo
[1,2-*c*]imidazole;

2-[4-[2-(2,3-Dimethylphenoxy)ethylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]
imidazole;

2-[4-[(Chroman-2-yl)methylamino]butyl]-1,4-dioxoperhydropyrido[1,2-*a*]pyrazine;

2-[4-[(Chroman-2-yl)methylamino]but-2-enyl]-1,4-dioxoperhydropyrrolo[1,2-*c*]
imidazole;

3-[4-[2-(*o*-Ethoxyphenoxy)ethylamino]butyl]-2,4-dioxothiazolidine;

3-[6-[2-(*o*-Ethoxyphenoxy)ethylamino]hexyl]-2,4-dioxothiazolidine;

3-[8-[2-(*o*-Ethoxyphenoxy)ethylamino]octyl]-2,4-dioxothiazolidine;

2-[4-[2-(*o*-Ethoxyphenoxy)ethylamino]butyl]-1,3-dioxoperhydroimidazo[1,5-*a*]pyridine;

2-[6-[2-(*o*-Ethoxyphenoxy)ethylamino]hexyl]-1,3-dioxoperhydroimidazo[1,5-*a*]pyridine;

2-[4-[(2-Quinolyl)methylamino]butyl]-1,3-dioxoperhydroimidazo[1,5-*a*]pyridine; and

2-[6-[(2-Quinolyl)methylamino]hexyl]-1,3-dioxoperhydroimidazo[1,5-*a*]pyridine.

23. **(new)** The use according to claim 21, wherein the isomer of the compound is
selected from the group consisting of:

2-[4-[(Chroman-2(*S*)-yl)methylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]imidazole;

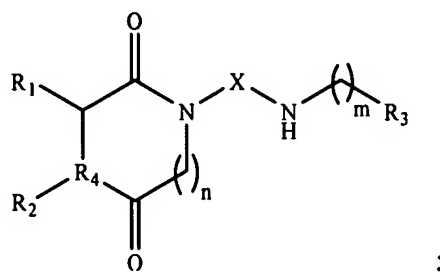
(*E*)-2-[4-[(Chroman-2-yl)methylamino]but-2-enyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]
imidazole; and

(*Z*)-2-[4-[(Chroman-2-yl)methylamino]but-2-enyl]-1,4-dioxoperhydropyrrolo[1,2-*c*]
imidazole.

24. **(new)** A method for preventing and/or treating a condition selected from the group
consisting of cerebral damage caused by thromboembolic stroke or traumatic brain damage,
Parkinson's disease, depression, migraine, pain, psychosis, mood disorder, and urinary tract
disorder in a subject in need of such prevention and/or treatment, wherein:

the method comprises administering to the subject a compound, a stereochemical isomer
of the compound, or a hydrate, solvate, or pharmaceutically acceptable salt of the compound or
isomer, wherein:

the compound corresponds in structure to formula I:



R₄ is selected from the group consisting of N and S;

if R₄ is S, then R₁ is H, and R₂ is absent;

if R₄ is N, then R₁ and R₂ are H or are methylene groups bound together forming with the heterocyclic ring a 5- or 6- membered ring;

n is zero or 1;

X is selected from the group consisting of C₂-C₁₀-alkylene, C₂-C₁₀-alkenylyne, and -CH₂-Y-CH₂-;

Y is phenyl;

m is 1 or 2;

R₃ is selected from the group consisting of chroman-2-yl, 2-quinolyl, and phenoxy, wherein:

the quinolyl, the aromatic ring of the chromanyl, and the phenyl ring of the phenoxy are optionally substituted with one or more substituents independently selected from the group consisting of C₁-C₆-alkoxy, C₁-C₆-alkyl, halogen, C₂-C₆-alkenyl, halo-(C₁-C₆)-alkyl, halo-(C₁-C₆)-alkoxy, phenyl, phenyl(C₁-C₆)-alkyl, phenoxy, C₁-C₆-alkylcarbonyl, phenylcarbonyl, phenyl(C₁-C₆)-alkylcarbonyl, C₁-C₆-alkoxycarbonyl, phenyl(C₁-C₆)-alkoxycarbonyl, C₁-C₆-alkyl-carbonylamino, hydroxy, cyano, nitro, amino, N-(C₁-C₆)-alkylamino, N,N-(C₁-C₆)-dialkylamino, carboxy, sulfo, sulfamoyl, sulfonylamino, (C₁-C₆)-alkylaminosulfonyl, and (C₁-C₆)-alkylsulfonylamino, wherein:

the C₁-C₆-alkyl portion of any of the alkyl-comprising substituents is optionally substituted with a substituent independently selected from the group consisting of hydroxy and amino; or

the phenyl ring of the phenoxy is substituted by two neighbouring residues, which together with the phenyl ring to which they are attached form tetrahydronaphthyl;

provided that the compound is not

2-[4-[(chroman-2-yl)methylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-c]imidazole,
3-[4-[(chroman-2-yl)methylamino]butyl]-2,4-dioxothiazolidine,
3-[5-[(chroman-2-yl)methylamino]pentyl]-2,4-dioxothiazolidine,
3-[6-[(chroman-2-yl)methylamino]hexyl]-2,4-dioxothiazolidine,
2-[4-[2-(phenoxy)ethylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-c]imidazole, or
3-[4-[2-(phenoxy)ethylamino]butyl]-2,4-dioxothiazolidine.

25. **(new)** The method according to claim 24, wherein the subject is human.

26. **(new)** The method according to claim 24, wherein the condition comprises migraine.

27. **(new)** The method according to claim 24, wherein the condition comprises pain.

28. **(new)** The method according to claim 24, wherein the condition comprises cerebral damage caused by thromboembolic stroke or traumatic brain damage.

29. **(new)** The method according to claim 24, wherein the condition comprises Parkinson's disease.

30. **(new)** The method according to claim 24, wherein the condition comprises depression.

31. **(new)** The method according to claim 24, wherein the condition comprises psychosis.

32. **(new)** The method according to claim 31, wherein the psychosis comprises schizophrenia.

33. **(new)** The method according to claim 24, wherein the condition comprises a mood disorder.

34. **(new)** The method according to claim 33, wherein the mood disorder comprises an anxiety disorder.

35. **(new)** The method according to claim 33, wherein the mood disorder comprises an aggressive disorder.

36. **(new)** The method according to claim 24, wherein the condition comprises a urinary tract disorder.

37. **(new)** The method according to claim 36, wherein the urinary tract disorder comprises urinary incontinence.

38. **(new)** The method according to claim 24, wherein an effective amount of the compound, isomer, hydrate, solvate, or salt is administered to the subject.

39. **(new)** The method according to claim 24, wherein the compound is selected from the group consisting of:

- 2-[4-[(Chroman-2-yl)methylamino]butyl]-1,3-dioxoperhydroimidazo[1,5-*a*]pyridine;
- 2-[4-[(Chroman-2-yl)methylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-*a*]pyrazine;
- 2-[5-[(Chroman-2-yl)methylamino]pentyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]imidazole;
- 2-[6-[(Chroman-2-yl)methylamino]hexyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]imidazole;
- 2-[3-[(Chroman-2-yl)methylamino]propyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]imidazole;
- 3-[8-[(Chroman-2-yl)methylamino]octyl]-2,4-dioxothiazolidine;
- 2-[4-[(Chroman-2-yl)methylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]imidazole;
- 2-[8-[(Chroman-2-yl)methylamino]octyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]imidazole;
- 2-[3-[[[(Chroman-2-yl)methylamino]methyl]benzyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]imidazole;
- 2-[4-[[[(Chroman-2-yl)methylamino]methyl]benzyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]imidazole;
- 2-[4-[(Chroman-2-yl)methylamino]but-2-enyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]imidazole;
- 2-[4-[2-(*o*-Methoxyphenoxy)ethylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]imidazole;
- 2-[4-[2-(*m*-Methoxyphenoxy)ethylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]

imidazole;

2-[4-[2-(*o*-Bromophenoxy)ethylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]imidazole;
2-[4-[2-(*m*-Bromophenoxy)ethylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]imidazole;
2-[4-[2-(*o*-Ethylphenoxy)ethylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]imidazole;
2-[4-[2-(*m*-Ethylphenoxy)ethylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]imidazole;
2-[4-[2-(*o*-Isopropylphenoxy)ethylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]

imidazole;

2-[4-[(2-quinolyl)methylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]imidazole;
2-[4-[2-(*o*-Ethoxyphenoxy)ethylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]imidazole;
2-[4-[2-(*o*-Isopropoxyphenoxy)ethylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]

imidazole;

2-[4-[2-[*m*-(Trifluoromethyl)phenoxy]ethylamino]butyl]-1,3-dioxoperhydropyrrolo
[1,2-*c*]imidazole;

2-[4-[2-(1,1'-Biphenyl-2-yloxy)ethylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]

imidazole;

2-[4-[2-[*o*-(Acetylamino)phenoxy]ethylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]

imidazole;

2-[4-[2-[*m*-(Acetylamino)phenoxy]ethylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]

imidazole;

2-[4-[2-[*o*-(Ethoxycarbonyl)phenoxy]ethylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]

imidazole;

2-[4-[2-(5,6,7,8-Tetrahydronaphth-1-yloxy)ethylamino]butyl]-1,3-dioxoperhydropyrrolo
[1,2-*c*]imidazole;

2-[4-[2-(2,3-Dimethylphenoxy)ethylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]

imidazole;

2-[4-[(Chroman-2-yl)methylamino]butyl]-1,4-dioxoperhydropyrrolo[1,2-*a*]pyrazine;

2-[4-[(Chroman-2-yl)methylamino]but-2-enyl]-1,4-dioxoperhydropyrrolo[1,2-*c*]

imidazole;

3-[4-[2-(*o*-Ethoxyphenoxy)ethylamino]butyl]-2,4-dioxothiazolidine;

3-[6-[2-(*o*-Ethoxyphenoxy)ethylamino]hexyl]-2,4-dioxothiazolidine;
3-[8-[2-(*o*-Ethoxyphenoxy)ethylamino]octyl]-2,4-dioxothiazolidine;
2-[4-[2-(*o*-Ethoxyphenoxy)ethylamino]butyl]-1,3-dioxoperhydroimidazo[1,5-*a*]pyridine;
2-[6-[2-(*o*-Ethoxyphenoxy)ethylamino]hexyl]-1,3-dioxoperhydroimidazo[1,5-*a*]pyridine;
2-[4-[(2-Quinolyl)methylamino]butyl]-1,3-dioxoperhydroimidazo[1,5-*a*]pyridine; and
2-[6-[(2-Quinolyl)methylamino]hexyl]-1,3-dioxoperhydroimidazo[1,5-*a*]pyridine.

40. **(new)** The method according to claim 24, wherein the isomer of the compound is selected from the group consisting of:

2-[4-[(Chroman-2(S)-yl)methylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]imidazole;
(*E*)-2-[4-[(Chroman-2-yl)methylamino]but-2-enyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]
imidazole; and
(*Z*)-2-[4-[(Chroman-2-yl)methylamino]but-2-enyl]-1,4-dioxoperhydropyrrolo[1,2-*c*]
imidazole.

41. **(new)** The compound, isomer, hydrate, solvate, or salt according to claim 1, wherein the isomer of the compound is selected from the group consisting of:

2-[4-[(Chroman-2(S)-yl)methylamino]butyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]imidazole;
(*E*)-2-[4-[(Chroman-2-yl)methylamino]but-2-enyl]-1,3-dioxoperhydropyrrolo[1,2-*c*]
imidazole; and
(*Z*)-2-[4-[(Chroman-2-yl)methylamino]but-2-enyl]-1,4-dioxoperhydropyrrolo[1,2-*c*]
imidazole.